



Full wwPDB X-ray Structure Validation Report ⓘ

Nov 25, 2017 – 07:43 PM EST

PDB ID : 5MI5
Title : BtGH84 mutant with covalent modification by MA3 in complex with PUGNAc
Authors : Darby, J.F.; Davies, G.J.; Hubbard, R.E.
Deposited on : unknown
Resolution : 2.15 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20030345
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : rb-20030345

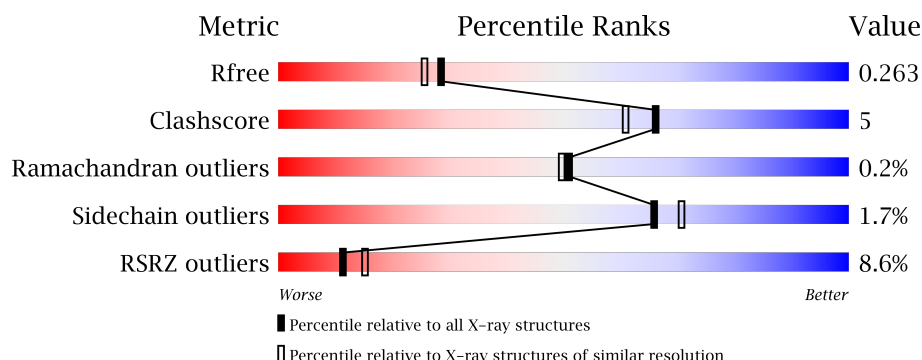
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.15 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	100719	1170 (2.16-2.16)
Clashscore	112137	1278 (2.16-2.16)
Ramachandran outliers	110173	1256 (2.16-2.16)
Sidechain outliers	110143	1255 (2.16-2.16)
RSRZ outliers	101464	1175 (2.16-2.16)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	727	

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 5786 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

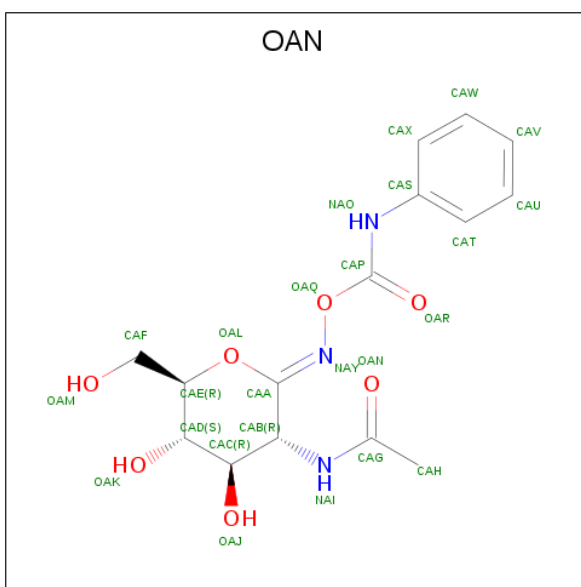
- Molecule 1 is a protein called O-GlcNAcase BT_4395.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	677	5490	3510	931	1030	19	0	1	0

There are 15 discrepancies between the modelled and reference sequences:

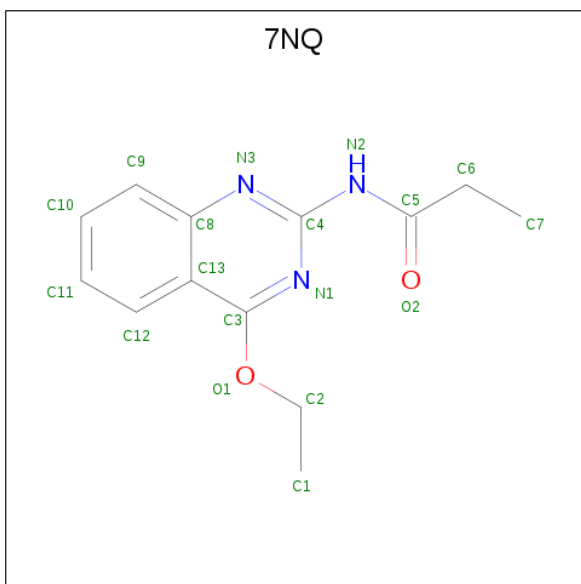
Chain	Residue	Modelled	Actual	Comment	Reference
A	-10	MET	-	initiating methionine	UNP Q89ZI2
A	-9	GLY	-	expression tag	UNP Q89ZI2
A	-8	SER	-	expression tag	UNP Q89ZI2
A	-7	SER	-	expression tag	UNP Q89ZI2
A	-6	HIS	-	expression tag	UNP Q89ZI2
A	-5	HIS	-	expression tag	UNP Q89ZI2
A	-4	HIS	-	expression tag	UNP Q89ZI2
A	-3	HIS	-	expression tag	UNP Q89ZI2
A	-2	HIS	-	expression tag	UNP Q89ZI2
A	-1	HIS	-	expression tag	UNP Q89ZI2
A	0	GLN	-	expression tag	UNP Q89ZI2
A	1	TRP	-	expression tag	UNP Q89ZI2
A	420	SER	CYS	engineered mutation	UNP Q89ZI2
A	550	CYS	TYR	engineered mutation	UNP Q89ZI2
A	654	SER	CYS	engineered mutation	UNP Q89ZI2

- Molecule 2 is O-(2-ACETAMIDO-2-DEOXY D-GLUCOPYRANOSYLIDENE) AMINO-N-PHENYLCARBAMATE (three-letter code: OAN) (formula: C₁₅H₁₉N₃O₇).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			25	15	3	7		

- Molecule 3 is {N}-(4-ethoxyquinazolin-2-yl)propanamide (three-letter code: 7NQ) (formula: $C_{13}H_{15}N_3O_2$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			18	13	3	2		

- Molecule 4 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total 1	Ca 1	0	0

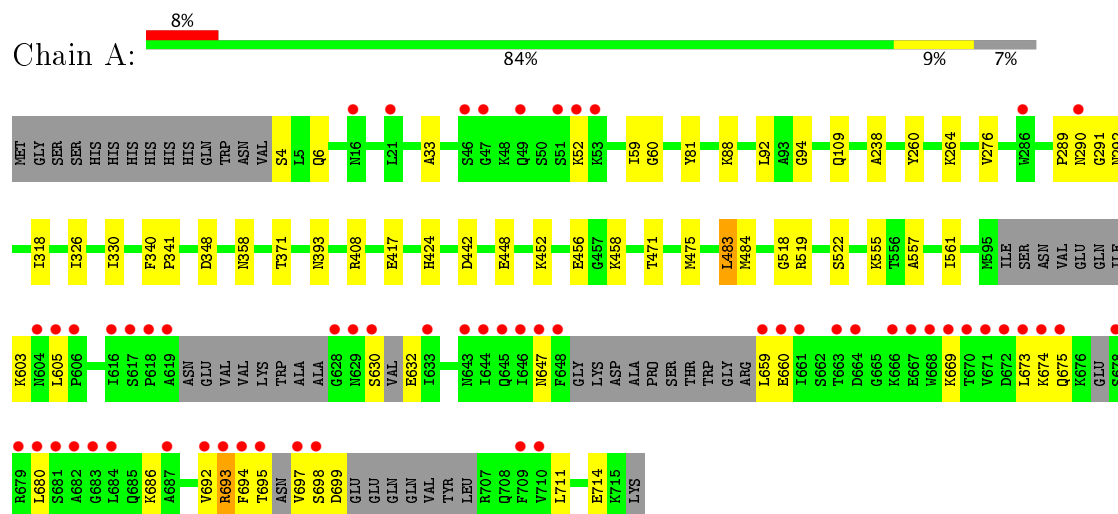
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	252	Total 252	O 252	0	0

3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

• Molecule 1: O-GlcNAcase BT_4395



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	198.22Å 52.12Å 112.78Å 90.00° 113.24° 90.00°	Depositor
Resolution (Å)	51.81 – 2.15 51.81 – 2.15	Depositor EDS
% Data completeness (in resolution range)	97.2 (51.81-2.15) 99.8 (51.81-2.15)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.92 (at 2.16Å)	Xtriage
Refinement program	PHENIX (1.11.1 _2575: ???)	Depositor
R, R_{free}	0.233 , 0.272 0.221 , 0.263	Depositor DCC
R_{free} test set	2903 reflections (5.02%)	DCC
Wilson B-factor (Å ²)	28.7	Xtriage
Anisotropy	0.620	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 47.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	5786	wwPDB-VP
Average B, all atoms (Å ²)	49.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.99% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: OAN, CA, 7NQ

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.28	0/5623	0.47	1/7607 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	693	ARG	NE-CZ-NH2	-8.69	115.95	120.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	5490	0	5412	50	0
2	A	25	0	19	0	0
3	A	18	0	0	0	0
4	A	1	0	0	0	0
5	A	252	0	0	8	1
All	All	5786	0	5431	50	1

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 5.

All (50) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:448:GLU:HB3	1:A:452:LYS:HZ3	0.98	1.08
1:A:448:GLU:HB3	1:A:452:LYS:NZ	1.71	1.05
1:A:448:GLU:O	1:A:452:LYS:HD3	1.64	0.96
1:A:442:ASP:OD2	5:A:901:HOH:O	1.87	0.93
1:A:483:LEU:HD23	1:A:484:MET:CE	2.06	0.84
1:A:448:GLU:CB	1:A:452:LYS:HZ3	1.88	0.83
1:A:483:LEU:HD23	1:A:484:MET:HE2	1.63	0.80
1:A:6:GLN:H	1:A:109:GLN:HE22	1.30	0.78
1:A:448:GLU:CB	1:A:452:LYS:NZ	2.44	0.78
1:A:318:ILE:H	1:A:358:ASN:HD22	1.33	0.76
1:A:693:ARG:NH2	5:A:902:HOH:O	2.11	0.75
1:A:660:GLU:OE2	1:A:693:ARG:NH2	2.23	0.72
1:A:603:LYS:HD3	1:A:605:LEU:HD23	1.72	0.70
1:A:673:LEU:HD23	1:A:680:LEU:HB3	1.80	0.62
1:A:318:ILE:H	1:A:358:ASN:ND2	1.98	0.61
1:A:483:LEU:HD23	1:A:484:MET:HE3	1.81	0.60
1:A:698:SER:OG	1:A:699:ASP:N	2.34	0.60
1:A:260:TYR:CZ	1:A:264:LYS:HD3	2.38	0.59
1:A:693:ARG:HD3	1:A:695:THR:OG1	2.02	0.59
1:A:555:LYS:NZ	5:A:914:HOH:O	2.36	0.58
1:A:52:LYS:O	1:A:52:LYS:NZ	2.24	0.57
1:A:260:TYR:CE1	1:A:264:LYS:HD3	2.40	0.56
1:A:659:LEU:HD12	5:A:902:HOH:O	2.04	0.56
1:A:693:ARG:HD2	1:A:694:PHE:N	2.22	0.55
1:A:660:GLU:OE2	1:A:693:ARG:CZ	2.59	0.50
1:A:424:HIS:HE1	5:A:901:HOH:O	1.95	0.50
1:A:660:GLU:OE2	1:A:693:ARG:NH1	2.45	0.49
1:A:4:SER:N	5:A:923:HOH:O	2.45	0.48
1:A:417:GLU:N	1:A:417:GLU:OE2	2.37	0.48
1:A:289:PRO:O	5:A:903:HOH:O	2.20	0.47
1:A:291:GLY:O	5:A:903:HOH:O	2.20	0.47
1:A:33:ALA:HB2	1:A:60:GLY:HA2	1.96	0.46
1:A:630:SER:O	1:A:632:GLU:N	2.49	0.46
1:A:471:THR:O	1:A:475:MET:HG3	2.15	0.46
1:A:326:ILE:O	1:A:330:ILE:HG12	2.15	0.46
1:A:557:ALA:HB1	1:A:561:ILE:HB	1.99	0.45
1:A:456:GLU:HG3	1:A:458:LYS:HD2	1.99	0.45
1:A:603:LYS:HB3	1:A:603:LYS:HE2	1.76	0.44
1:A:519:ARG:HG2	1:A:519:ARG:H	1.59	0.44
1:A:695:THR:HG22	1:A:697:VAL:N	2.33	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:238:ALA:HA	1:A:276:VAL:O	2.18	0.43
1:A:81:TYR:HA	1:A:94:GLY:HA2	2.01	0.42
1:A:59:ILE:HA	1:A:92:LEU:O	2.20	0.42
1:A:692:VAL:HG21	1:A:711:LEU:HD22	2.01	0.41
1:A:340:PHE:CD1	1:A:341:PRO:HA	2.55	0.41
1:A:292:ASN:OD1	1:A:292:ASN:N	2.54	0.41
1:A:686:LYS:NZ	1:A:714:GLU:OE1	2.44	0.41
1:A:442:ASP:N	1:A:442:ASP:OD2	2.54	0.41
1:A:340:PHE:O	1:A:371:THR:OG1	2.31	0.40
1:A:348:ASP:N	1:A:348:ASP:OD1	2.53	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:A:959:HOH:O	5:A:959:HOH:O[2_555]	1.90	0.30

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	A	662/727 (91%)	630 (95%)	31 (5%)	1 (0%)	51 50

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	518	GLY

5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	597/640 (93%)	587 (98%)	10 (2%)	66	71

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	88	LYS
1	A	290	ASN
1	A	393	ASN
1	A	408	ARG
1	A	483	LEU
1	A	522	SER
1	A	647	ASN
1	A	669	LYS
1	A	674	LYS
1	A	675	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	109	GLN
1	A	198	ASN
1	A	358	ASN
1	A	393	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates

There are no carbohydrates in this entry.

5.6 Ligand geometry

Of 3 ligands modelled in this entry, 1 is monoatomic - leaving 2 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	OAN	A	801	-	23,26,26	1.48	2 (8%)	23,35,35	1.76	5 (21%)
3	7NQ	A	802	1	19,19,19	2.57	5 (26%)	23,25,25	3.11	8 (34%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	OAN	A	801	-	-	2/12/35/35	0/2/2/2
3	7NQ	A	802	1	-	0/9/9/9	0/2/2/2

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	801	OAN	OAQ-NAY	-4.34	1.33	1.44
3	A	802	7NQ	C13-C8	-3.89	1.36	1.42
2	A	801	OAN	OAL-CAE	-3.68	1.41	1.46
3	A	802	7NQ	O2-C5	-2.01	1.19	1.23
3	A	802	7NQ	C5-N2	4.12	1.45	1.35
3	A	802	7NQ	C4-N2	5.02	1.45	1.38
3	A	802	7NQ	O1-C3	7.50	1.40	1.35

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	802	7NQ	C4-N2-C5	-8.97	116.40	130.28
3	A	802	7NQ	C13-C3-N1	-7.02	119.50	124.45
2	A	801	OAN	CAS-NAO-CAP	-5.25	117.70	126.41
2	A	801	OAN	CAC-CAD-CAE	-4.06	103.07	110.22
3	A	802	7NQ	C13-C8-N3	-3.68	119.04	122.84
3	A	802	7NQ	N3-C4-N1	-3.03	121.64	126.23
3	A	802	7NQ	C2-O1-C3	-3.00	114.51	117.58
2	A	801	OAN	OAK-CAD-CAC	-2.48	104.97	110.36
3	A	802	7NQ	O2-C5-N2	-2.02	119.87	123.67
2	A	801	OAN	CAA-CAB-NAI	2.07	112.59	109.26
2	A	801	OAN	CAX-CAS-CAT	2.25	122.17	119.04
3	A	802	7NQ	C4-N1-C3	3.04	120.54	115.18
3	A	802	7NQ	O1-C3-C13	5.71	119.93	114.45

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	801	OAN	NAY-OAQ-CAP-OAR
2	A	801	OAN	NAY-OAQ-CAP-NAO

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	677/727 (93%)	0.46	58 (8%) 11 15	24, 43, 94, 123	0

All (58) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	47	GLY	6.0
1	A	659	LEU	5.4
1	A	619	ALA	5.0
1	A	604	ASN	5.0
1	A	51	SER	4.5
1	A	694	PHE	4.5
1	A	290	ASN	4.4
1	A	630	SER	4.3
1	A	605	LEU	4.2
1	A	671	VAL	4.2
1	A	693	ARG	3.8
1	A	681	SER	3.8
1	A	661	ILE	3.8
1	A	668	TRP	3.6
1	A	666	LYS	3.5
1	A	682	ALA	3.5
1	A	679	ARG	3.5
1	A	53	LYS	3.5
1	A	698	SER	3.4
1	A	710	VAL	3.3
1	A	628	GLY	3.3
1	A	673	LEU	3.3
1	A	646	ILE	3.1
1	A	692	VAL	3.1
1	A	695	THR	3.1
1	A	633	ILE	3.1
1	A	616	ILE	3.0

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Mol	Chain	Res	Type	RSRZ
1	A	21	LEU	3.0
1	A	680	LEU	3.0
1	A	660	GLU	3.0
1	A	644	ILE	2.9
1	A	52	LYS	2.9
1	A	678	SER	2.9
1	A	648	PHE	2.8
1	A	683	GLY	2.7
1	A	618	PRO	2.7
1	A	286	TRP	2.6
1	A	674	LYS	2.6
1	A	670	THR	2.6
1	A	647	ASN	2.6
1	A	669	LYS	2.4
1	A	687	ALA	2.4
1	A	667	GLU	2.4
1	A	697	VAL	2.4
1	A	709	PHE	2.4
1	A	46	SER	2.4
1	A	643	ASN	2.4
1	A	606	PRO	2.3
1	A	629	ASN	2.3
1	A	16	ASN	2.3
1	A	49	GLN	2.3
1	A	672	ASP	2.2
1	A	684	LEU	2.2
1	A	645	GLN	2.1
1	A	617	SER	2.1
1	A	664	ASP	2.0
1	A	675	GLN	2.0
1	A	663	THR	2.0

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
3	7NQ	A	802	18/18	0.91	0.14	1.56	36,46,60,62	0
2	OAN	A	801	25/25	0.89	0.15	0.96	29,44,72,73	0
4	CA	A	803	1/1	0.95	0.22	-	57,57,57,57	0

6.5 Other polymers [i](#)

There are no such residues in this entry.